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 NEWS  1           Web Page URLs for STN Seminar Schedule - N. America
 NEWS  2  Sep 29   The Philippines Inventory of Chemicals and Chemical
                   Substances (PICCS) has been added to CHEMLIST
 NEWS  3  Oct 27   New Extraction Code PAX now available in Derwent
                   Files
 NEWS  4  Oct 27   SET ABBREVIATIONS and SET PLURALS extended in
                   Derwent World Patents Index files
 NEWS  5  Oct 27   Patent Assignee Code Dictionary now available
                   in Derwent Patent Files
 NEWS  6  Oct 27   Plasdoc Key Serials Dictionary and Echoing added to
                   Derwent Subscriber Files WPIDS and WPIX
 NEWS  7  Nov 29   Derwent announces further increase in updates for DWPI
 NEWS  8  Dec  5   French Multi-Disciplinary Database PASCAL Now on STN
 NEWS  9  Dec  5   Trademarks on STN - New DEMAS and EUMAS Files
 NEWS 10  Dec 15   2001 STN Pricing
 NEWS 11  Dec 17   Merged CEABA-VTB for chemical engineering and
                   biotechnology
 NEWS 12  Dec 17   Corrosion Abstracts on STN
 NEWS 13  Dec 17   SYNTHLINE from Prous Science now available on STN
 NEWS 14  Dec 17   The CA Lexicon available in the CAPLUS and CA files
 NEWS 15  Jan 05   AIDSLINE is being removed from STN
 NEWS 16  Feb 06   Engineering Information Encompass files have new names
 NEWS 17  Feb 16   TOXLINE no longer being updated

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 NEWS HOURS    STN Operating Hours Plus Help Desk Availability
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Enter NEWS followed by the item number or name to see news on that
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FILE 'HOME' ENTERED AT 17:11:24 ON 12 MAR 2001

=> file reg
```

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.15	0.15

FILE 'REGISTRY' ENTERED AT 17:11:32 ON 12 MAR 2001

STRUCTURE FILE UPDATES: 11 MAR 2001 HIGHEST RN 326792-71-6
DICTIONARY FILE UPDATES: 11 MAR 2001 HIGHEST RN 326792-71-6

TSCA INFORMATION NOW CURRENT THROUGH July 8, 2000

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

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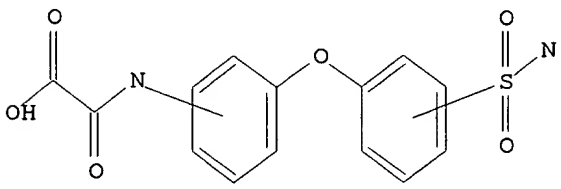
Uploading 533219.str

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS
L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 17:12:11 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 11 TO ITERATE

100.0% PROCESSED 11 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:	ONLINE	**COMPLETE**
	BATCH	**COMPLETE**
PROJECTED ITERATIONS:	22 TO	418
PROJECTED ANSWERS:	1 TO	80

L2 1 SEA SSS SAM L1

=> s l1 ful

FULL SEARCH INITIATED 17:12:17 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 213 TO ITERATE

100.0% PROCESSED 213 ITERATIONS 33 ANSWERS
SEARCH TIME: 00.00.01

L3 33 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	133.56	133.71

FILE 'CAPLUS' ENTERED AT 17:12:24 ON 12 MAR 2001
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FILE COVERS 1967 - 12 Mar 2001 VOL 134 ISS 12
FILE LAST UPDATED: 9 Mar 2001 (20010309/ED)

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=> s l3

L4 2 L3

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=> d k4 1- ibib abs hitstr

'K4' IS NOT A VALID FORMAT FOR FILE 'CAPLUS'

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ABS --------- GI and AB
ALL --------- BIB, AB, IND, RE
APPS -------- AI, PRAI
BIB --------- AN, plus Bibliographic Data and PI table (default)
CAN --------- List of CA abstract numbers without answer numbers
CBIB -------- AN, plus Compressed Bibliographic Data
DALL -------- ALL, delimited (end of each field identified)
DMAX -------- MAX, delimited for post-processing
FAM --------- AN, PI and PRAI in table, plus Patent Family data
FBIB -------- AN, BIB, plus Patent FAM
IND --------- Indexing data
IPC --------- International Patent Classifications
MAX --------- ALL, plus Patent FAM, RE
PATS -------- PI, SO
SAM --------- CC, SX, TI, ST, IT
SCAN -------- CC, SX, TI, ST, IT (random display, no answer numbers;
              SCAN must be entered on the same line as the DISPLAY,
              e.g., D SCAN or DISPLAY SCAN)
STD --------- BIB, IPC, and NCL

IABS -------- ABS, indented with text labels
IALL -------- ALL, indented with text labels
IBIB -------- BIB, indented with text labels
IMAX -------- MAX, indented with text labels
ISTD -------- STD, indented with text labels

OBIB -------- AN, plus Bibliographic Data (original)
OIBIB ------- OBIB, indented with text labels

SBIB -------- BIB, no citations
SIBIB ------- IBIB, no citations

HIT --------- Fields containing hit terms
HITIND ------ IC, ICA, ICI, NCL, CC and index field (ST and IT)
              containing hit terms
HITRN ------- HIT RN and its text modification
HITSTR ------ HIT RN, its text modification, its CA index name, and
              its structure diagram
FHITSTR ----- First HIT RN, its text modification, its CA index name, and
              its structure diagram
KWIC -------- Hit term plus 20 words on either side
OCC --------- Number of occurrence of hit term and field in which it occurs

To display a particular field or fields, enter the display field
codes.  For a list of the display field codes, enter HELP DFIELDS at
an arrow prompt (=>).  Examples of formats include: TI; TI,AU; BIB,ST;
TI,IND; TI,SO.  You may specify the format fields in any order and the
information will be displayed in the same order as the format
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ENTER DISPLAY FORMAT (BIB):ibib abs hitstr

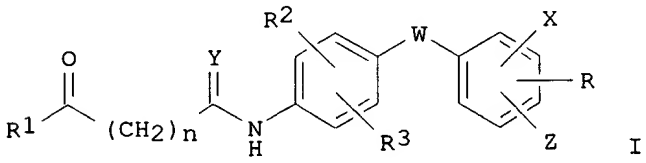
YOU HAVE REQUESTED DATA FROM 2 ANSWERS - CONTINUE? Y/(N):y

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 2000:707138 CAPLUS
DOCUMENT NUMBER: 133:266609
TITLE: Preparation of (4-phenoxyphenyl)oxamic acid derivatives and analogs as hypolipidemics
INVENTOR(S): Kukkola, Paivi Jaana
PATENT ASSIGNEE(S): Novartis A.-G., Switz.; Novartis-Erfindungen
SOURCE: PCT Int. Appl., 53 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000058279	A1	20001005	WO 2000-EP2683	20000327

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

PRIORITY APPLN. INFO.: US 1999-280105 19990329
OTHER SOURCE(S): MARPAT 133:266609
GI



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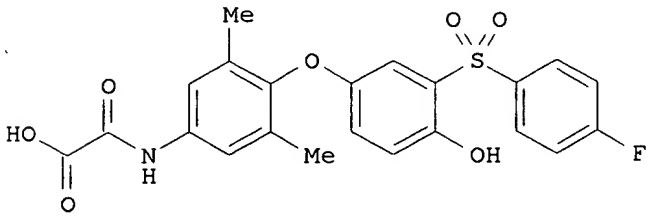


II

AB The title compds. (I) [wherein W = O, S, S(O) or SO2; X = SR4, S(O)R4, SO2R4, SO2NR5R6, or CONR5R6; Y = O or H2; Z = H, halogen, OH, or (un)substituted (ar)alkoxy, acyloxy, or alkoxycarbonyloxy; R = H, halogen, CF3, or (cyclo)alkyl; R1 = OH, (un)substituted (cyclo)alkoxy, (hetero)aryloxy, or (hetero)aralkoxy, or -NR5R6; R2 = H, halogen, or alkyl; R3 = halogen or alkyl; R4 is (un)substituted (ar)alkyl, (hetero)aryl, or heteroaralkyl; R5, R6, and R7 = independently H, (un)substituted (cyclo)alkyl, (hetero)aryl, or (hetero)aralkyl; or R5 and R6 combined = alkylene optionally interrupted by O, S, S(O), SO2, or NR7 which together with the nitrogen atom to which they are attached form a 5- to 7-membered ring; n = 0-4] were prepd. I demonstrated potent binding to the triiodothyronine (T3) nuclear receptor, which is indicative of upregulation of LDL receptor activity and enhancement of the clearance of LDL-cholesterol from the circulation. I also reduced lipoprotein (a) levels and are useful for the treatment and prevention of occlusive cardiovascular conditions implicated by Lp(a). For example, 2-(4-fluorobenzensulfonyl)benzene-1,4-diol (prepn. given) was coupled with 4-chloro-3,5-dimethylnitrobenzene in the presence of NaH, and the product reduced using Pd/C. Amidation with di-Et oxalate, followed by deesterification, gave II. In an in vitro T3 nuclear receptor binding assay using Sprague-Dawley rat liver nuclei and plasma membrane prepns., II gave an IC50 of 0.17 nM. II significantly lowered serum cholesterol at a daily dose of about 20 .mu.g/kg p.o. in male Sprague-Dawley rats and about 10 .mu.g/kg p.o. in normocholesterolemic dogs. Lp(a) levels in normolipemic cynomolgus monkeys were lowered by about 40% after a 4 wk treatment with II at a daily oral dose of 75 .mu.g/kg. Thus, I are useful in the prevention and treatment of diseases assocd. with an imbalance of thyroid hormones, such as hypo- and hyperthyroidism, obesity, osteoporosis, and depression, and for lowering LDL cholesterol and Lp(a) levels.

IT **298694-79-8P**, N-[4-[3-(2,2-Dimethylpropylsulfamoyl)-4-hydroxyphenoxy]-3,5-dimethylphenyl]oxamic acid **298694-80-1P**, N-[4-(4-Hydroxy-3-phenylsulfamoylphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-81-2P**, N-[4-[3-(4-Fluorophenylsulfamoyl)-4-hydroxyphenoxy]-3,5-dimethylphenyl]oxamic acid **298694-82-3P**, N-[4-[3-(2-Fluorophenylsulfamoyl)-4-hydroxyphenoxy]-3,5-dimethylphenyl]oxamic acid **298694-84-5P**, N-[4-[3-(3-Fluorophenylsulfamoyl)-4-hydroxyphenoxy]-3,5-dimethylphenyl]oxamic acid **298694-85-6P**, N-[4-[4-Hydroxy-3-(4-methoxyphenylsulfamoyl)phenoxy]-3,5-dimethylphenyl]oxamic acid **298694-86-7P**, N-[4-[3-(4-Fluorobenzylsulfamoyl)-4-hydroxyphenoxy]-3,5-dimethylphenyl]oxamic acid **298694-87-8P**, N-[4-[4-Hydroxy-3-(N-methyl-N-phenylsulfamoyl)phenoxy]-3,5-dimethylphenyl]oxamic acid **298694-88-9P**, N-[4-(4-Hydroxy-3-propylsulfamoylphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-89-0P**, N-[4-(4-Hydroxy-3-isopropylsulfamoylphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-90-3P**, N-[4-(3-Butylsulfamoyl-4-hydroxyphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-91-4P**, N-[4-(4-Hydroxy-3-

isobutylsulfamoylphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-92-5P**, N-[4-(3-t-Butylsulfamoyl-4-hydroxyphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-93-6P**, N-[4-(3-Cyclohexylsulfamoyl-4-hydroxyphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-94-7P**, N-[4-(3-Dimethylsulfamoyl-4-hydroxyphenoxy)-3,5-dimethylphenyl]oxamic acid **298694-97-0P**, N-[4-[4-Hydroxy-3-(2-methoxyethylsulfamoyl)phenoxy]-3,5-dimethylphenyl]oxamic acid **298695-00-8P**, N-[4-[4-Hydroxy-3-(pyridin-3-ylsulfamoyl)phenoxy]-3,5-dimethylphenyl]oxamic acid **298695-01-9P**, N-[4-[4-Hydroxy-3-(1-methyl-6-oxo-1,6-dihydropyridin-3-ylsulfamoyl)phenoxy]-3,5-dimethylphenyl]oxamic acid **298695-03-1P**, N-[4-[3-(4-Fluorophenylsulfamoyl)phenoxy]-3,5-dimethylphenyl]oxamic acid

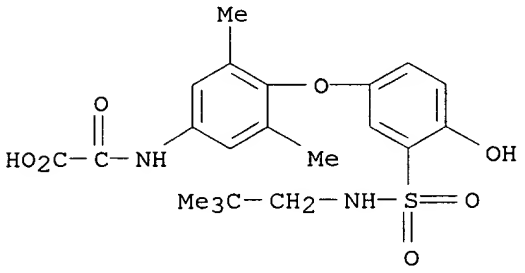
298695-04-2P, N-[4-[3-(4-Fluorophenylsulfamoyl)-4-hydroxyphenoxy]-3-methylphenyl]oxamic acid

RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of (4-phenoxyphenyl)oxamic acid derivs. and analogs as hypolipidemics by coupling phenols with 4-chloronitrobenzenes, redn. to the amines, and amidation with oxalates)

RN 298694-79-8 CAPLUS

CN Acetic acid, [[4-[3-[[(2,2-dimethylpropyl)amino]sulfonyl]-4-hydroxyphenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 298694-80-1 CAPLUS

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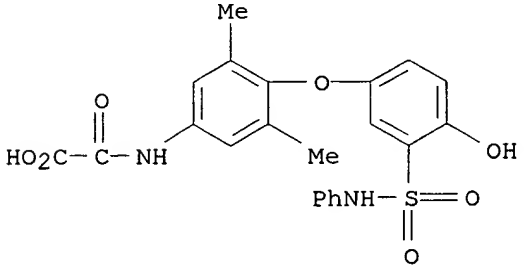

RN 298694-81-2 CAPLUS
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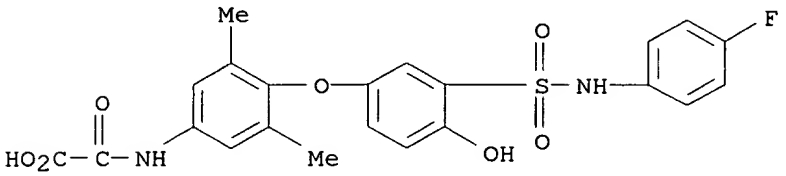

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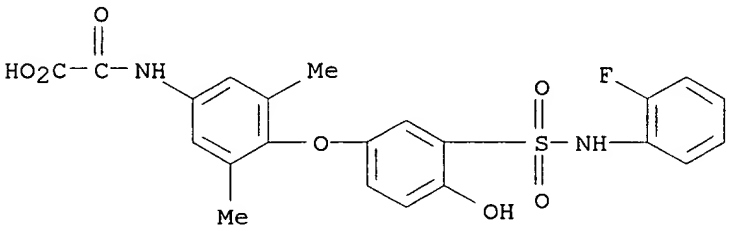

RN 298694-84-5 CAPLUS
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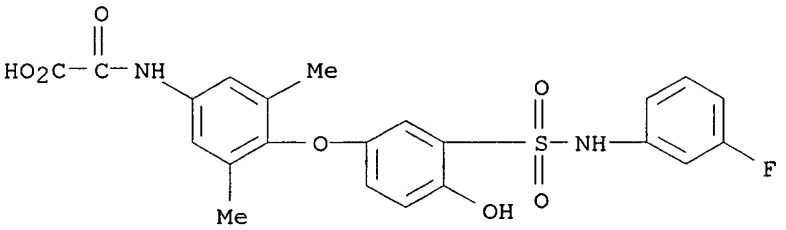

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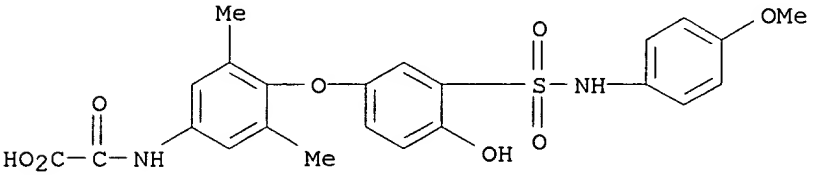

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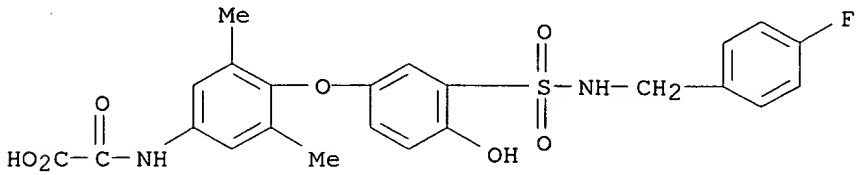

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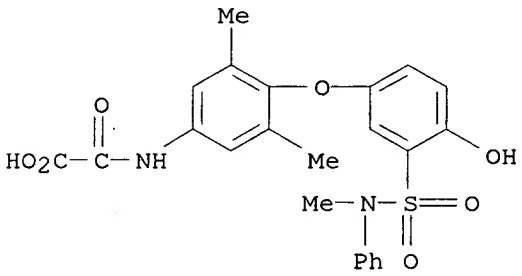

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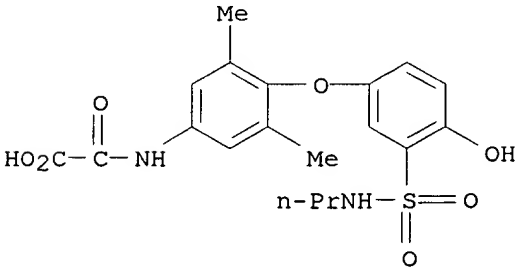

RN 298694-89-0 CAPLUS
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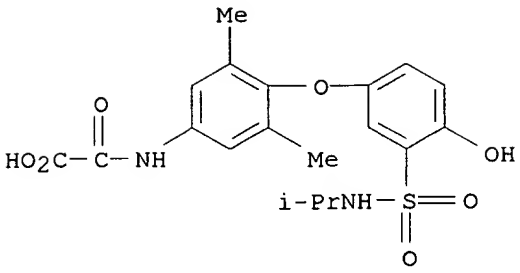

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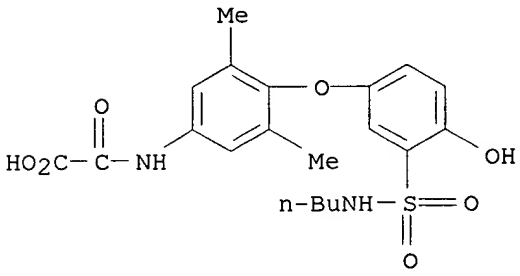

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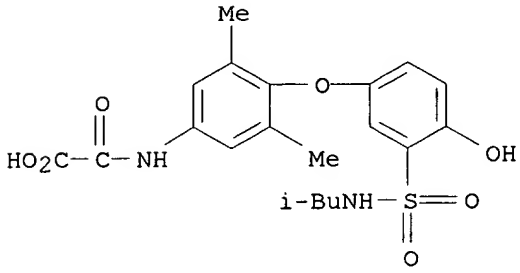

RN 298694-92-5 CAPLUS
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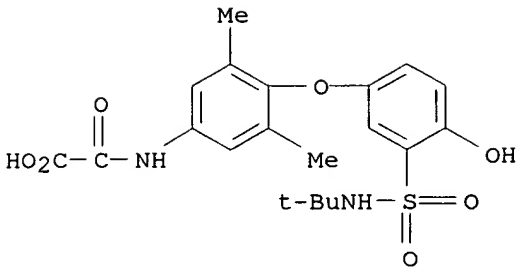

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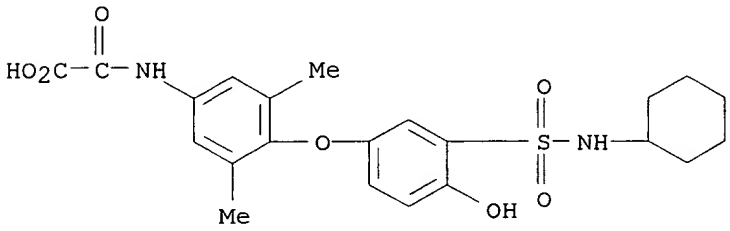

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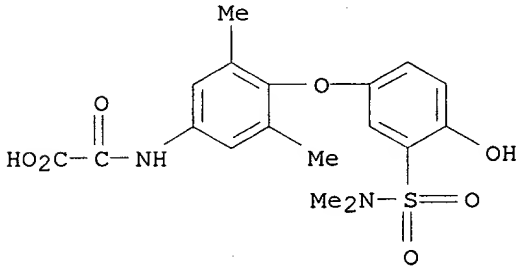

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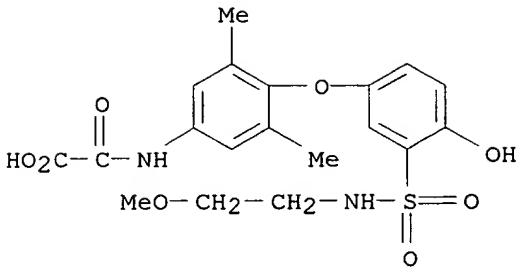
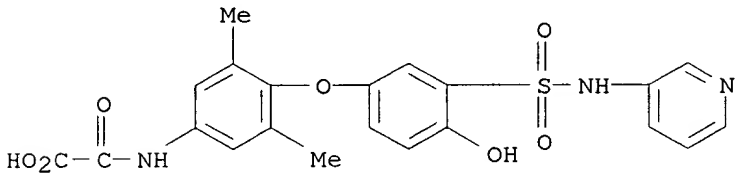
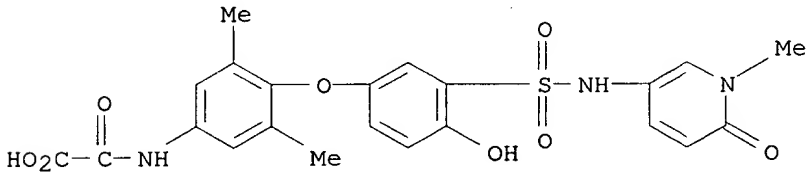


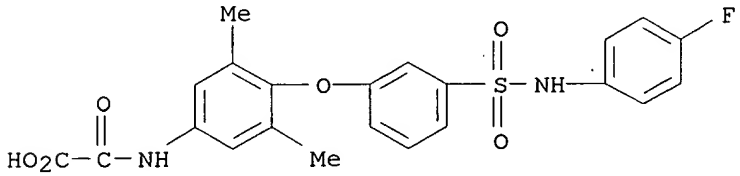
RN 298695-00-8 CAPLUS
CN Acetic acid, [[4-[4-hydroxy-3-[(3-pyridinylamino)sulfonyl]phenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 298695-01-9 CAPLUS
CN Acetic acid, [[4-[3-[[(1,6-dihydro-1-methyl-6-oxo-3-pyridinyl)amino]sulfonyl]-4-hydroxyphenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 298695-03-1 CAPLUS
CN Acetic acid, [[4-[3-[[(4-fluorophenyl)amino]sulfonyl]phenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 298695-04-2 CAPLUS
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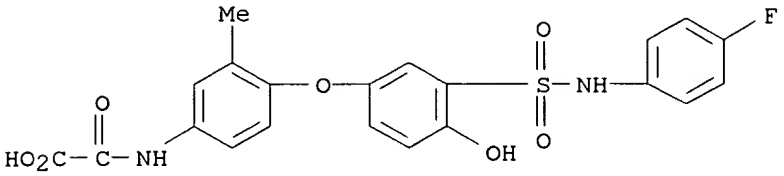
methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



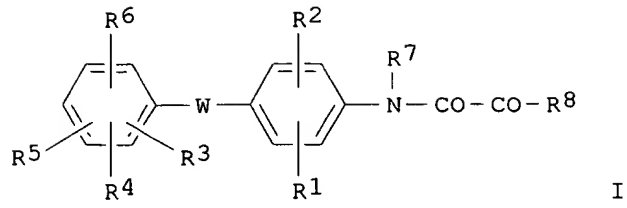
REFERENCE COUNT: 2
REFERENCE(S): (1) Ciba Geigy Ag; EP 0580550 A 1994 CAPLUS
(2) Yokoyama, N; JOURNAL OF MEDICINAL CHEMISTRY 1995, V38, P695 CAPLUS

L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 2000:628106 CAPLUS
DOCUMENT NUMBER: 133:207681
TITLE: Preparation of 4-(sulfamoylphenoxy)phenyloxamic acids and derivatives as thyroid receptor ligands
INVENTOR(S): Chiang, Yuan-Ching Phoebe; Dow, Robert Lee
PATENT ASSIGNEE(S): Pfizer Products Inc., USA
SOURCE: PCT Int. Appl., 128 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000051971	A1	20000908	WO 2000-IB183	20000221

W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

PRIORITY APPLN. INFO.: US 1999-122292 19990301
OTHER SOURCE(S): MARPAT 133:207681
GI



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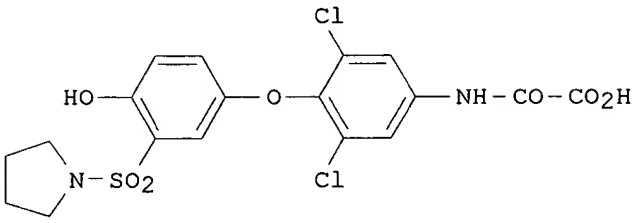


II

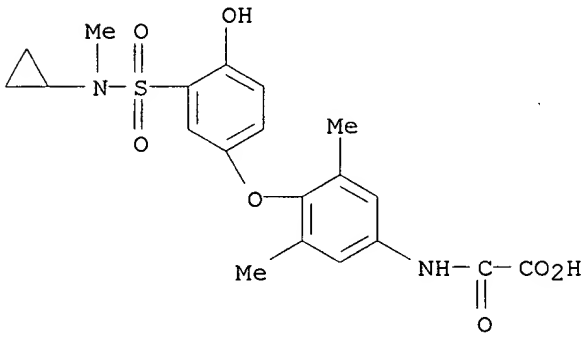
AB The title compds. (I) [wherein R1-R3 = independently H, halo, alkyl, CF3, CN, OCF3, or alkoxy; R4 = H or (un)substituted alkyl; or R3 and R4 together form an (un)substituted carbocyclic ring, (CH2)b, or a heterocyclic ring, Q(CH2)c or (CH2)jQ(CH2)k; b = 3-7; c = 2-6; j and k = independently 2-6; Q = O, S, or NR1; R5 = F, OH, alkoxy, or carboxy; or R4 and R5 together form a heterocyclic ring; R6 = H, halo, alkyl, or CF3; R7 = H or alkyl; R8 = OH, alkoxy, or (un)substituted amino; W = O, S(O)d, CH2, NH, or N(alkyl); d = 0-2], prodrugs, geometric and optical isomers, and pharmaceutically acceptable salts were prepd. as thyroid receptor ligands. Thus, 2',6'-dichloro-4-methoxy-4'-nitrodiphenyl ether was treated with ClSO2H and pyrrolidine in two steps to give 1-[5-(2,6-dichloro-4-nitrophenoxy)-2-methoxybenzenesufonyl]pyrrolidine. Demethylation using BCl3, followed by redn. using Pd/C, addn. of di-Et oxalate, and deesterification, yielded II. An in vivo oxygen consumption assay designed to evaluate the efficacy and cardiac effects of tissue-selective thyroid hormone agonists and a thyroid hormone receptor (TR.alpha. and TR.beta.) binding assay for thyromimetic compds. are described (no data). I are useful for the treatment of obesity, hyperlipidemia, glaucoma, cardiac arrhythmia, skin disorders, thyroid disease, hypothyroidism, and related disorders and diseases, such as diabetes mellitus, atherosclerosis, hypertension, coronary heart disease, hypercholesteremia, depression, and osteoporosis. An anorectic agent or lipase inhibitor may be administered with I to treat these conditions.

IT **290349-36-9P 290349-37-0P 290349-38-1P,**

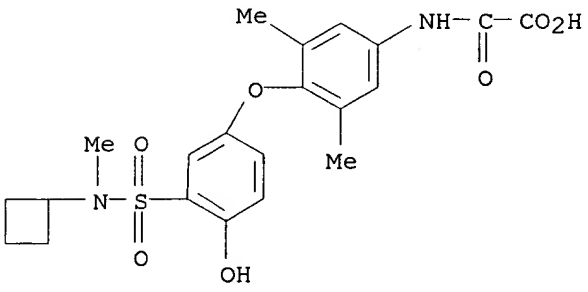
N-[4-(3-(Cyclopropylsulfamoyl)-4-hydroxyphenoxy)-3,5-dimethylphenyl]oxamic acid **290349-39-2P**, N-[4-(3-(Cyclobutylsulfamoyl)-4-hydroxyphenoxy)-3,5-dimethylphenyl]oxamic acid **290349-63-2P**, N-[3-Chloro-4-(3-cyclopropylsulfamoyl-4-hydroxyphenoxy)-5-methylphenyl]oxamic acid **290349-64-3P**, N-[3-Chloro-4-(3-cyclobutylsulfamoyl-4-hydroxyphenoxy)-5-methylphenyl]oxamic acid **290349-65-4P**, N-[3-Chloro-4-(3-cyclopentylsulfamoyl-4-hydroxyphenoxy)-5-methylphenyl]oxamic acid **290349-66-5P**, N-[3-Chloro-4-(3-cyclohexylsulfamoyl-4-hydroxyphenoxy)-5-

methylphenyl]oxamic acid **290349-67-6P**, N-[3-Chloro-4-(4-hydroxy-3-sulfamoylphenoxy)-5-methylphenyl]oxamic acid **290349-68-7P**, N-[3-Chloro-4-[3-(4-fluorophenylsulfamoyl)-4-hydroxyphenoxy]-5-methylphenyl]oxamic acid **290349-69-8P**, N-[3-Chloro-4-(4-hydroxy-3-propylsulfamoylphenoxy)-5-methylphenyl]oxamic acid **290349-70-1P**, N-[4-(3-Butylsulfamoyl-4-hydroxyphenoxy)-3-chloro-5-methylphenyl]oxamic acid **290349-78-9P**, N-[3,5-Dichloro-4-(4-hydroxy-3-isopropylsulfamoylphenoxy)phenyl]oxamic acid
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of 4-(sulfamoylphenoxy)phenyloxamic acids and derivs. as thyroid receptor ligands by treatment of 4-methoxy-4'-nitrodiphenyl ethers with ClSO3H and amines, redn., and amidation with oxalates)

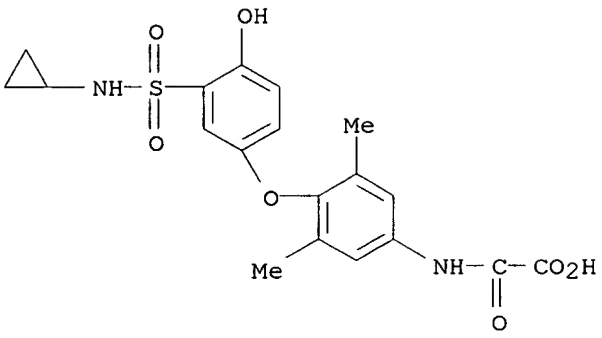
RN 290349-36-9 CAPLUS
CN Acetic acid, [[4-[3-[(cyclopropylmethylamino)sulfonyl]-4-hydroxyphenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



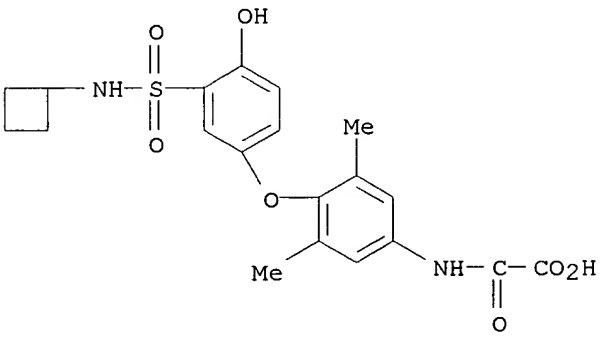
RN 290349-37-0 CAPLUS
CN Acetic acid, [[4-[3-[(cyclobutylmethylamino)sulfonyl]-4-hydroxyphenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



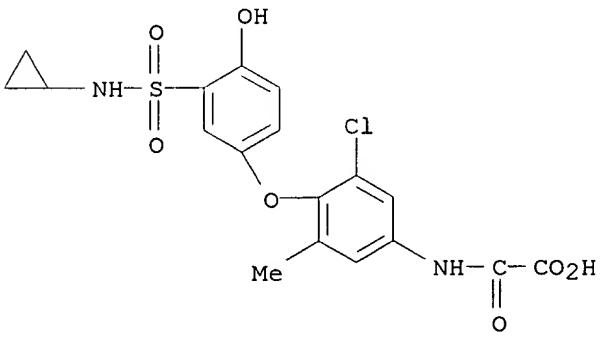
RN 290349-38-1 CAPLUS
CN Acetic acid, [[4-[3-[(cyclopropylamino)sulfonyl]-4-hydroxyphenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 290349-39-2 CAPLUS
CN Acetic acid, [[4-[3-[(cyclobutylamino)sulfonyl]-4-hydroxyphenoxy]-3,5-dimethylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 290349-63-2 CAPLUS
CN Acetic acid, [[3-chloro-4-[3-[(cyclopropylamino)sulfonyl]-4-hydroxyphenoxy]-5-methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 290349-64-3 CAPLUS
CN Acetic acid,
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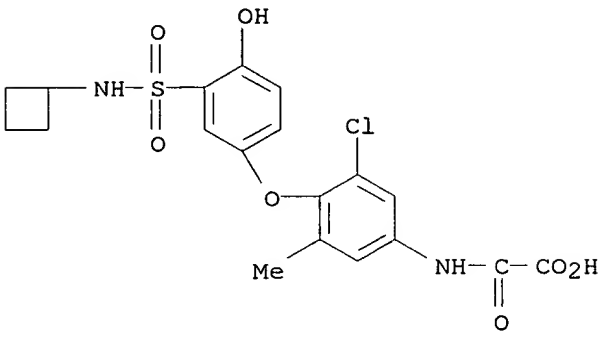
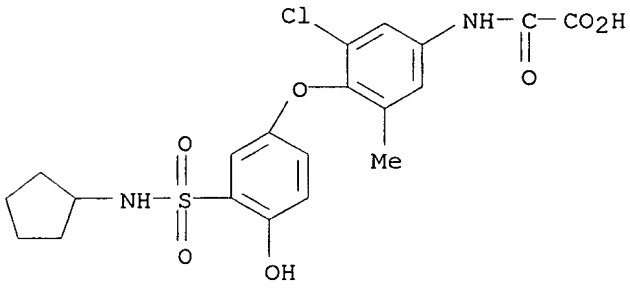
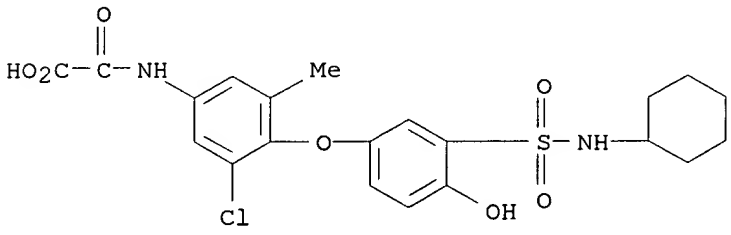
5-methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



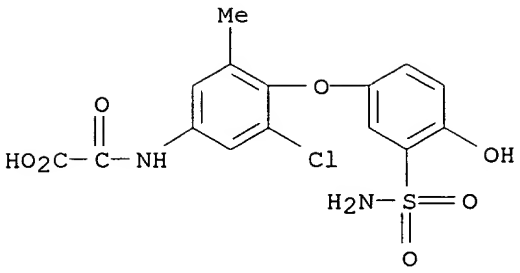
RN 290349-65-4 CAPLUS
CN Acetic acid, [[3-chloro-4-[3-[(cyclopentylamino)sulfonyl]-4-hydroxyphenoxy]-5-methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



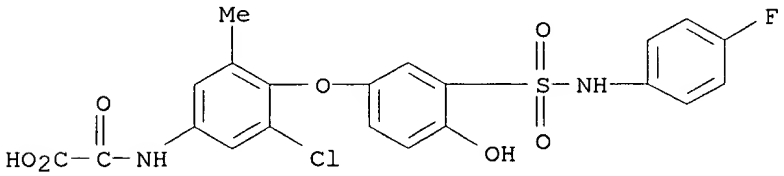
RN 290349-66-5 CAPLUS
CN Acetic acid, [[3-chloro-4-[3-[(cyclohexylamino)sulfonyl]-4-hydroxyphenoxy]-5-methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 290349-67-6 CAPLUS
CN Acetic acid, [[4-[3-(aminosulfonyl)-4-hydroxyphenoxy]-3-chloro-5-methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 290349-68-7 CAPLUS
CN Acetic acid, [[3-chloro-4-[3-[[(4-fluorophenyl)amino]sulfonyl]-4-hydroxyphenoxy]-5-methylphenyl]amino]oxo- (9CI) (CA INDEX NAME)



RN 290349-69-8 CAPLUS
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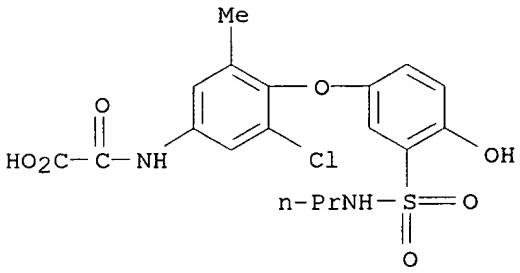

RN 290349-70-1 CAPLUS
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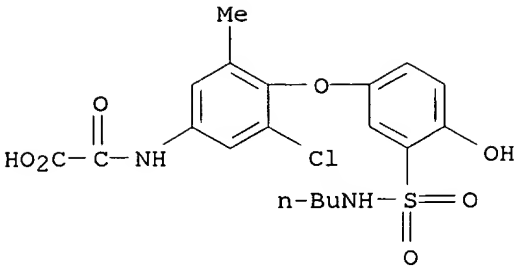
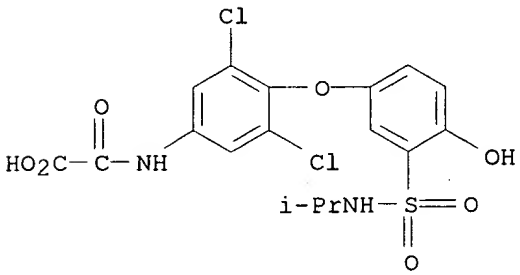


RN 290349-78-9 CAPLUS
CN Acetic acid, [[3,5-dichloro-4-[4-hydroxy-3-[[(1-methylethyl)amino]sulfonyl]phenoxy]phenyl]amino]oxo- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 4
REFERENCE(S):
(1) Apelqvist, T; WO 0007972 A 2000 CAPLUS
(2) Ciba Geigy Ag; EP 0580550 A 1994 CAPLUS
(3) Taylor, A; MOLECULAR PHARMACOLOGY 1997, V52(3), P542 CAPLUS
(4) Yokoyama, N; JOURNAL OF MEDICINAL CHEMISTRY 1995, V38, P695 CAPLUS

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